

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051921\
 Data File : BP005606.D
 Acq On : 19 May 2021 11:07
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad

5/20/2021 11:35:31 AM

Quant Time: May 19 13:20:07 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 17 15:35:47 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.658	152	10224	20.00	ng	# 0.00
21) Naphthalene-d8	10.446	136	35731	20.00	ng	# 0.00
39) Acenaphthene-d10	14.310	164	31496	20.00	ng	0.00
64) Phenanthrene-d10	17.069	188	79525	20.00	ng	# 0.00
76) Chrysene-d12	21.175	240	113130	20.00	ng	# 0.00
86) Perylene-d12	23.469	264	121718	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.252	112	41354	77.62	ng	0.00
7) Phenol-d6	6.858	99	53510	80.55	ng	0.00
23) Nitrobenzene-d5	8.828	82	70870	78.07	ng	0.00
42) 2,4,6-Tribromophenol	15.816	330	64671	72.36	ng	0.00
45) 2-Fluorobiphenyl	12.928	172	205391	73.56	ng	0.00
79) Terphenyl-d14	19.645	244	531873	79.84	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.140	88	8156	39.18	ng	# 88
3) Pyridine	3.558	79	19320	40.96	ng	# 75
4) n-Nitrosodimethylamine	3.470	42	15219	40.12	ng	# 4
6) Aniline	6.999	93	28552	39.82	ng	# 69
8) 2-Chlorophenol	7.228	128	23085	40.34	ng	91
9) Benzaldehyde	6.816	77	15721	47.03	ng	# 71
10) Phenol	6.887	94	25920	39.01	ng	# 45
11) bis(2-Chloroethyl)ether	7.099	93	17850	38.90	ng	91
12) 1,3-Dichlorobenzene	7.546	146	28151	36.98	ng	# 80
13) 1,4-Dichlorobenzene	7.693	146	29238	38.53	ng	97
14) 1,2-Dichlorobenzene	8.005	146	28585	39.31	ng	95
15) Benzyl Alcohol	7.922	79	25147	39.24	ng	# 81
16) 2,2'-oxybis(1-Chloropr...	8.199	45	8442	37.26	ng	# 52
17) 2-Methylphenol	8.128	107	17975	39.39	ng	# 75
18) Hexachloroethane	8.716	117	12816	40.78	ng	87
19) n-Nitroso-di-n-propyla...	8.475	70	20722	42.60	ng	89
20) 3+4-Methylphenols	8.463	107	25247	41.62	ng	88
22) Acetophenone	8.493	105	37312	37.02	ng	# 84
24) Nitrobenzene	8.869	77	35495	38.98	ng	# 88
25) Isophorone	9.399	82	54250	38.59	ng	# 95
26) 2-Nitrophenol	9.581	139	13657	35.82	ng	# 75
27) 2,4-Dimethylphenol	9.652	122	19190	38.02	ng	# 83
28) bis(2-Chloroethoxy)met...	9.881	93	21745	36.95	ng	98
29) 2,4-Dichlorophenol	10.122	162	30302	38.23	ng	97
30) 1,2,4-Trichlorobenzene	10.316	180	39093	38.61	ng	99
31) Naphthalene	10.499	128	72568	37.77	ng	98
32) Benzoic acid	9.863	122	13652	35.01	ng	# 77
33) 4-Chloroaniline	10.634	127	28231	37.72	ng	# 89
34) Hexachlorobutadiene	10.769	225	35177	41.66	ng	94
35) Caprolactam	11.452	113	6643m	36.43	ng	
36) 4-Chloro-3-methylphenol	11.781	107	28317	40.08	ng	94
37) 2-Methylnaphthalene	12.122	142	59904	38.87	ng	95
38) 1-Methylnaphthalene	12.340	142	57263	39.36	ng	95
40) 1,2,4,5-Tetrachloroben...	12.493	216	53503	36.97	ng	99
41) Hexachlorocyclopentadiene	12.457	237	17800	34.25	ng	98
43) 2,4,6-Trichlorophenol	12.751	196	32278	35.70	ng	98

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44) 2,4,5-Trichlorophenol	12.840	196	33827	34.67	ng	92
46) 1,1'-Biphenyl	13.146	154	88402	36.45	ng	96
47) 2-Chloronaphthalene	13.187	162	74054	36.68	ng	98
48) 2-Nitroaniline	13.416	65	21969	38.30	ng	96
49) Acenaphthylene	14.034	152	104818	36.03	ng	99
50) Dimethylphthalate	13.787	163	97792	35.97	ng	98
51) 2,6-Dinitrotoluene	13.916	165	21542	35.28	ng	87
52) Acenaphthene	14.375	154	65645	36.27	ng	97
53) 3-Nitroaniline	14.251	138	15659	35.80	ng	93
54) 2,4-Dinitrophenol	14.481	184	13449	36.52	ng	# 83
55) Dibenzofuran	14.716	168	126252	37.09	ng	94
56) 4-Nitrophenol	14.598	139	13525	40.96	ng	# 54
57) 2,4-Dinitrotoluene	14.710	165	32695	36.32	ng	# 82
58) Fluorene	15.363	166	104271	37.51	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.957	232	37134m	38.06	ng	
60) Diethylphthalate	15.151	149	96598	36.68	ng	96
61) 4-Chlorophenyl-phenyle...	15.363	204	68275	39.25	ng	97
62) 4-Nitroaniline	15.422	138	14831	35.29	ng	# 66
63) Azobenzene	15.657	77	106781	39.34	ng	86
65) 4,6-Dinitro-2-methylph...	15.481	198	22151	35.15	ng	95
66) n-Nitrosodiphenylamine	15.587	169	88133	37.97	ng	94
67) 4-Bromophenyl-phenylether	16.257	248	48848	39.35	ng	98
68) Hexachlorobenzene	16.369	284	61393	39.00	ng	99
69) Atrazine	16.545	200	41303	38.10	ng	98
70) Pentachlorophenol	16.734	266	34194	40.94	ng	95
71) Phenanthrene	17.110	178	172821	38.24	ng	98
72) Anthracene	17.204	178	169034	37.69	ng	98
73) Carbazole	17.486	167	146392	37.07	ng	100
74) Di-n-butylphthalate	18.039	149	164307	37.90	ng	# 96
75) Fluoranthene	19.092	202	237682	38.24	ng	97
77) Benzidine	19.292	184	68591	39.17	ng	97
78) Pyrene	19.445	202	245165	38.28	ng	99
80) Butylbenzylphthalate	20.328	149	77766	37.58	ng	93
81) Benzo(a)anthracene	21.157	228	277904	37.91	ng	99
82) 3,3'-Dichlorobenzidine	21.098	252	100886	36.53	ng	# 97
83) Chrysene	21.210	228	270714	38.10	ng	99
84) Bis(2-ethylhexyl)phtha...	21.080	149	117274	37.06	ng	# 99
85) Di-n-octyl phthalate	21.969	149	199014	37.44	ng	# 91
87) Indeno(1,2,3-cd)pyrene	25.857	276	365438	35.69	ng	# 96
88) Benzo(b)fluoranthene	22.769	252	306274	37.76	ng	# 99
89) Benzo(k)fluoranthene	22.816	252	302724	38.17	ng	# 98
90) Benzo(a)pyrene	23.369	252	282952	38.06	ng	# 98
91) Dibenzo(a,h)anthracene	25.857	278	321903	35.63	ng	# 97
92) Benzo(g,h,i)perylene	26.586	276	308977	36.96	ng	# 94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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